



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 07:06 PM UTC

PDB ID : 1Z6A / pdb_00001z6a
Title : Sulfolobus solfataricus SWI2/SNF2 ATPase core domain
Authors : Duerr, H.; Koerner, C.; Mueller, M.; Hickmann, V.; Hopfner, K.P.
Deposited on : 2005-03-22
Resolution : 3.00 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	NOT EXECUTED
EDS	:	NOT EXECUTED
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

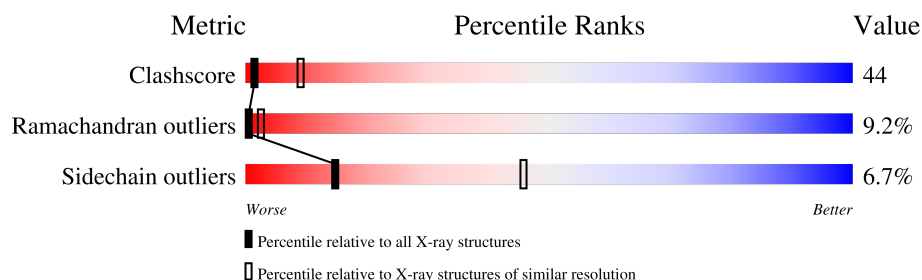
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.00 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
Clashscore	190562	2977 (3.00-3.00)
Ramachandran outliers	187476	2877 (3.00-3.00)
Sidechain outliers	187428	2880 (3.00-3.00)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	500	

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 3874 atoms, of which 0 are hydrogens and 0 are deuteriums.

In the tables below, the ZeroOcc column contains the number of atoms modelled with zero occupancy, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

- Molecule 1 is a protein called Helicase of the snf2/rad54 family.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	480	3831	2467	644	707	13	62	0	0

There are 24 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-12	MET	-	expression tag	UNP Q97XQ5
A	-11	GLY	-	expression tag	UNP Q97XQ5
A	-10	SER	-	expression tag	UNP Q97XQ5
A	-9	SER	-	expression tag	UNP Q97XQ5
A	-8	HIS	-	expression tag	UNP Q97XQ5
A	-7	HIS	-	expression tag	UNP Q97XQ5
A	-6	HIS	-	expression tag	UNP Q97XQ5
A	-5	HIS	-	expression tag	UNP Q97XQ5
A	-4	HIS	-	expression tag	UNP Q97XQ5
A	-3	HIS	-	expression tag	UNP Q97XQ5
A	-2	SER	-	expression tag	UNP Q97XQ5
A	-1	SER	-	expression tag	UNP Q97XQ5
A	0	GLY	-	expression tag	UNP Q97XQ5
A	1	LEU	-	expression tag	UNP Q97XQ5
A	2	VAL	-	expression tag	UNP Q97XQ5
A	3	PRO	-	expression tag	UNP Q97XQ5
A	4	ARG	-	expression tag	UNP Q97XQ5
A	5	GLY	-	expression tag	UNP Q97XQ5
A	6	SER	-	expression tag	UNP Q97XQ5
A	7	HIS	-	expression tag	UNP Q97XQ5
A	8	MET	-	expression tag	UNP Q97XQ5
A	9	ALA	-	expression tag	UNP Q97XQ5
A	11	SER	-	expression tag	UNP Q97XQ5
A	12	LYS	-	expression tag	UNP Q97XQ5

- Molecule 2 is MERCURY (II) ION (CCD ID: HG) (formula: Hg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	3	Total	Hg	0	0
			3	3		

- Molecule 3 is PHOSPHATE ION (CCD ID: PO4) (formula: O_4P).



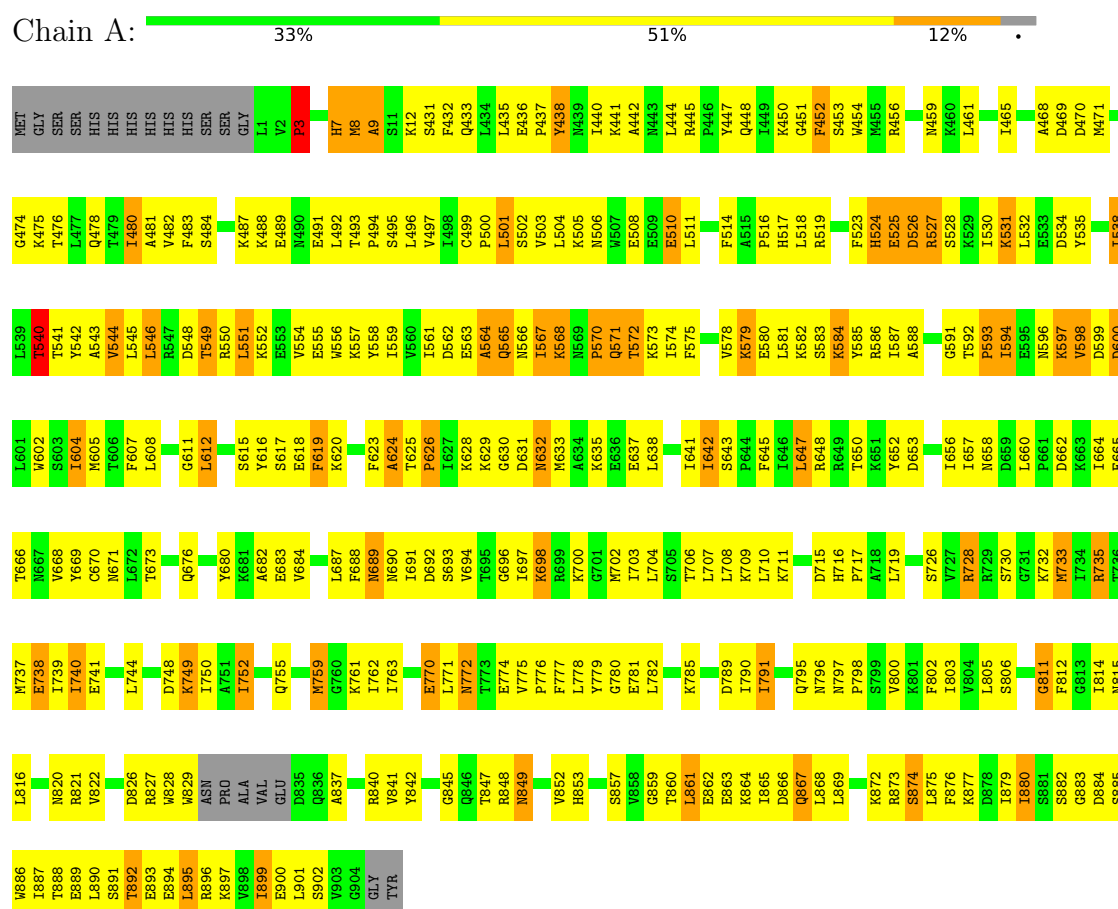
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	O	P	0	0
			5	4	1		
3	A	1	Total	O	P	0	0
			5	4	1		
3	A	1	Total	O	P	0	0
			5	4	1		
3	A	1	Total	O	P	0	0
			5	4	1		
3	A	1	Total	O	P	0	0
			5	4	1		
3	A	1	Total	O	P	0	0
			5	4	1		
3	A	1	Total	O	P	0	0
			5	4	1		

3 Residue-property plots

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

Note EDS was not executed.

- Molecule 1: Helicase of the *snf2/rad54* family



4 Data and refinement statistics

Xtriage (Phenix) and EDS were not executed - this section is therefore incomplete.

Property	Value	Source
Space group	P 41 21 2	Depositor
Cell constants a, b, c, α , β , γ	83.89Å 83.89Å 227.54Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	20.00 – 3.00	Depositor
% Data completeness (in resolution range)	(Not available) (20.00-3.00)	Depositor
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	CNS	Depositor
R, R_{free}	0.240 , 0.297	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	3874	wwPDB-VP
Average B, all atoms (Å ²)	91.0	wwPDB-VP

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: HG, PO4

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
1	A	0.52	0/3898	1.03	25/5252 (0.5%)

There are no bond length outliers.

All (25) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	584	LYS	N-CA-C	-8.14	104.50	114.75
1	A	531	LYS	N-CA-C	-6.82	102.01	112.99
1	A	759	MET	N-CA-C	-6.78	103.96	111.82
1	A	3	PRO	N-CA-CB	6.67	110.25	103.25
1	A	820	ASN	N-CA-C	-6.63	105.32	113.41
1	A	480	ILE	N-CA-C	-6.60	104.32	110.53
1	A	510	GLU	N-CA-C	6.24	119.37	111.69
1	A	849	ASN	N-CA-C	-6.17	102.39	110.53
1	A	658	ASN	N-CA-C	6.05	117.68	111.14
1	A	861	LEU	N-CA-C	-5.90	104.92	111.71
1	A	643	SER	CA-C-N	-5.82	113.66	120.12
1	A	643	SER	C-N-CA	-5.82	113.66	120.12
1	A	895	LEU	N-CA-C	-5.62	105.16	111.28
1	A	741	GLU	N-CA-C	5.57	117.43	111.36
1	A	874	SER	N-CA-C	-5.56	106.47	113.20
1	A	709	LYS	N-CA-C	-5.56	105.29	112.68
1	A	899	ILE	N-CA-C	5.41	116.19	110.72
1	A	875	LEU	N-CA-C	-5.38	107.73	114.56
1	A	540	THR	N-CA-C	5.29	116.56	108.42
1	A	475	LYS	N-CA-C	-5.23	105.76	111.82
1	A	796	ASN	N-CA-C	5.23	116.67	110.97
1	A	641	ILE	N-CA-C	5.20	115.92	110.62
1	A	698	LYS	N-CA-C	-5.14	107.17	112.93
1	A	749	LYS	N-CA-C	-5.04	100.08	108.34

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	600	ASP	N-CA-C	-5.03	106.32	112.90

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts ⓘ

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	3831	0	3926	336	1
2	A	3	0	0	0	0
3	A	40	0	0	2	0
All	All	3874	0	3926	336	1

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 44.

All (336) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:571:GLN:HG2	1:A:572:THR:H	1.11	1.07
1:A:568:LYS:O	1:A:570:PRO:HD3	1.53	1.07
1:A:650:THR:HG22	1:A:652:TYR:H	1.19	1.02
1:A:546:LEU:HD21	1:A:573:LYS:HE3	1.44	0.97
1:A:791:ILE:HD11	1:A:814:ILE:HG23	1.47	0.96
1:A:860:THR:HG22	1:A:862:GLU:H	1.29	0.95
1:A:650:THR:HG22	1:A:652:TYR:N	1.86	0.90
1:A:531:LYS:HG2	1:A:534:ASP:HB2	1.52	0.90
1:A:542:TYR:HB3	1:A:574:ILE:HD13	1.55	0.87
1:A:592:THR:O	1:A:594:ILE:HG13	1.73	0.87
1:A:491:GLU:C	1:A:557:LYS:HD3	2.02	0.84
1:A:571:GLN:HG2	1:A:572:THR:N	1.91	0.84
1:A:625:THR:OG1	1:A:626:PRO:HD3	1.77	0.83
1:A:519:ARG:HD2	1:A:535:TYR:HE2	1.43	0.81
1:A:748:ASP:OD2	1:A:821:ARG:NH1	2.13	0.81
1:A:883:GLY:HA2	1:A:886:TRP:HE1	1.45	0.81

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:650:THR:CG2	1:A:652:TYR:H	1.94	0.81
1:A:814:ILE:HD12	1:A:842:TYR:CE2	2.16	0.80
1:A:837:ALA:O	1:A:841:VAL:HG23	1.81	0.80
1:A:867:GLN:HG2	1:A:868:LEU:N	1.95	0.80
1:A:555:GLU:HG2	1:A:582:LYS:HE2	1.64	0.79
1:A:600:ASP:O	1:A:604:ILE:HG13	1.82	0.78
1:A:715:ASP:HA	1:A:732:LYS:HD2	1.64	0.78
1:A:454:TRP:CZ3	1:A:647:LEU:HB2	2.18	0.78
1:A:527:ARG:CZ	1:A:530:ILE:HG12	2.14	0.78
1:A:797:ASN:HD22	1:A:798:PRO:HD2	1.48	0.77
1:A:759:MET:HE3	1:A:762:ILE:HB	1.66	0.77
1:A:478:GLN:O	1:A:482:VAL:HG23	1.86	0.76
1:A:740:ILE:HD12	1:A:802:PHE:HE2	1.51	0.75
1:A:638:LEU:HG	1:A:642:ILE:HD11	1.67	0.75
1:A:592:THR:HG22	1:A:648:ARG:CZ	2.18	0.74
1:A:700:LYS:HE2	1:A:888:THR:OG1	1.87	0.74
1:A:504:LEU:HD21	1:A:508:GLU:OE1	1.88	0.74
1:A:495:SER:OG	1:A:558:TYR:HB2	1.87	0.74
1:A:638:LEU:O	1:A:642:ILE:HG13	1.86	0.74
1:A:593:PRO:HA	1:A:597:LYS:NZ	2.03	0.73
1:A:873:ARG:HG3	1:A:876:PHE:HD2	1.52	0.73
1:A:593:PRO:HA	1:A:597:LYS:HZ2	1.52	0.73
1:A:519:ARG:HD2	1:A:535:TYR:CE2	2.23	0.72
1:A:873:ARG:HA	1:A:876:PHE:HB3	1.71	0.72
1:A:863:GLU:O	1:A:866:ASP:HB3	1.91	0.71
1:A:631:ASP:OD1	1:A:633:MET:HB2	1.91	0.71
1:A:668:VAL:HG11	1:A:739:ILE:HD12	1.73	0.71
1:A:9:ALA:HA	1:A:432:PHE:CE1	2.27	0.70
1:A:873:ARG:HG2	1:A:877:LYS:HB2	1.73	0.70
1:A:869:LEU:HD21	1:A:886:TRP:CH2	2.27	0.70
1:A:563:GLU:O	1:A:564:ALA:O	2.10	0.69
1:A:597:LYS:NZ	1:A:597:LYS:HB3	2.07	0.69
1:A:504:LEU:O	1:A:508:GLU:HG3	1.93	0.69
1:A:676:GLN:OE1	1:A:730:SER:HA	1.92	0.69
1:A:465:ILE:HG22	1:A:645:PHE:HA	1.75	0.69
1:A:527:ARG:HD2	1:A:530:ILE:CD1	2.23	0.69
1:A:476:THR:O	1:A:480:ILE:HD13	1.94	0.68
1:A:605:MET:HE3	1:A:605:MET:HA	1.76	0.68
1:A:653:ASP:O	1:A:656:ILE:HG22	1.92	0.68
1:A:726:SER:HB2	1:A:728:ARG:HE	1.58	0.68
1:A:749:LYS:HE2	1:A:800:VAL:O	1.94	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:499:CYS:HB2	1:A:503:VAL:CG2	2.23	0.67
1:A:527:ARG:HD2	1:A:530:ILE:HD11	1.76	0.67
1:A:740:ILE:HD11	1:A:752:ILE:HD11	1.77	0.67
1:A:440:ILE:HD11	1:A:481:ALA:HA	1.74	0.66
1:A:688:PHE:HE1	1:A:896:ARG:HB2	1.59	0.66
1:A:444:LEU:HD12	1:A:444:LEU:H	1.60	0.65
1:A:883:GLY:O	1:A:887:ILE:HG13	1.96	0.65
1:A:556:TRP:O	1:A:583:SER:HA	1.95	0.65
1:A:561:ILE:HD11	1:A:578:VAL:HG11	1.78	0.64
1:A:728:ARG:CD	1:A:728:ARG:H	2.10	0.64
1:A:571:GLN:CG	1:A:572:THR:H	1.98	0.64
1:A:859:GLY:O	1:A:901:LEU:HD12	1.98	0.64
1:A:703:ILE:CG2	1:A:888:THR:HG22	2.28	0.63
1:A:441:LYS:N	1:A:441:LYS:HD2	2.12	0.63
1:A:688:PHE:O	1:A:691:ILE:HG22	1.97	0.63
1:A:688:PHE:C	1:A:690:ASN:H	2.07	0.62
1:A:733:MET:O	1:A:737:MET:HG2	1.99	0.62
1:A:527:ARG:HD3	1:A:550:ARG:CB	2.30	0.62
1:A:704:LEU:O	1:A:708:LEU:HG	2.00	0.61
1:A:735:ARG:HA	1:A:738:GLU:HG3	1.81	0.61
1:A:791:ILE:CD1	1:A:814:ILE:HA	2.30	0.61
1:A:860:THR:CG2	1:A:861:LEU:HD23	2.31	0.61
1:A:459:ASN:HB3	1:A:585:TYR:HD2	1.63	0.61
1:A:548:ASP:CG	1:A:549:THR:H	2.09	0.61
1:A:633:MET:HE2	1:A:633:MET:HA	1.83	0.60
1:A:543:ALA:C	1:A:545:LEU:H	2.09	0.60
1:A:471:MET:O	1:A:821:ARG:NH2	2.34	0.60
1:A:578:VAL:C	1:A:580:GLU:H	2.09	0.60
1:A:484:SER:O	1:A:487:LYS:HB3	2.02	0.60
1:A:555:GLU:HG2	1:A:582:LYS:CE	2.31	0.60
1:A:865:ILE:O	1:A:869:LEU:HG	2.02	0.60
1:A:523:PHE:HB3	1:A:540:THR:OG1	2.02	0.60
1:A:580:GLU:HG3	1:A:580:GLU:O	2.01	0.60
1:A:673:THR:H	1:A:676:GLN:NE2	2.00	0.59
1:A:527:ARG:HH21	1:A:550:ARG:CB	2.15	0.59
1:A:602:TRP:HB2	1:A:619:PHE:CD2	2.38	0.59
1:A:480:ILE:O	1:A:483:PHE:HB2	2.03	0.59
1:A:444:LEU:H	1:A:444:LEU:CD1	2.16	0.59
1:A:711:LYS:NZ	1:A:755:GLN:HE22	2.01	0.59
1:A:465:ILE:HG22	1:A:645:PHE:CA	2.33	0.58
1:A:891:SER:O	1:A:893:GLU:N	2.36	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:541:THR:HG22	1:A:542:TYR:N	2.17	0.58
1:A:715:ASP:OD2	1:A:732:LYS:HE3	2.03	0.58
1:A:564:ALA:C	1:A:566:ASN:H	2.10	0.58
1:A:444:LEU:HD12	1:A:444:LEU:N	2.17	0.58
1:A:688:PHE:HE2	1:A:899:ILE:HD11	1.67	0.58
1:A:887:ILE:O	1:A:890:LEU:HG	2.02	0.58
1:A:8:MET:O	1:A:9:ALA:HB2	2.03	0.58
1:A:857:SER:O	1:A:860:THR:HB	2.04	0.58
1:A:516:PRO:C	1:A:518:LEU:H	2.11	0.58
1:A:597:LYS:HB3	1:A:597:LYS:HZ3	1.67	0.58
1:A:620:LYS:HA	1:A:624:ALA:HB3	1.85	0.58
1:A:715:ASP:HB3	1:A:759:MET:SD	2.44	0.57
1:A:660:LEU:HD22	1:A:664:ILE:HD11	1.87	0.57
1:A:797:ASN:HD22	1:A:798:PRO:CD	2.17	0.57
1:A:592:THR:O	1:A:594:ILE:N	2.29	0.57
1:A:775:VAL:HG12	1:A:775:VAL:O	2.04	0.57
1:A:563:GLU:OE2	1:A:570:PRO:HB2	2.04	0.57
1:A:687:LEU:O	1:A:691:ILE:HB	2.04	0.57
1:A:717:PRO:C	1:A:719:LEU:N	2.61	0.57
1:A:827:ARG:HH22	1:A:866:ASP:HB2	1.68	0.57
1:A:899:ILE:HG13	1:A:900:GLU:N	2.19	0.57
1:A:849:ASN:HB2	3:A:961:PO4:O2	2.04	0.57
1:A:689:ASN:N	1:A:689:ASN:HD22	2.02	0.56
1:A:771:LEU:O	1:A:772:ASN:C	2.48	0.56
1:A:890:LEU:HB2	1:A:895:LEU:HD13	1.88	0.56
1:A:564:ALA:HB1	1:A:565:GLN:NE2	2.21	0.55
1:A:795:GLN:HE21	1:A:815:ASN:ND2	2.05	0.55
1:A:586:ARG:NH2	1:A:608:LEU:O	2.40	0.55
1:A:540:THR:HG23	1:A:544:VAL:HB	1.89	0.55
1:A:605:MET:HG3	1:A:642:ILE:HG21	1.89	0.55
1:A:504:LEU:C	1:A:504:LEU:HD23	2.32	0.55
1:A:445:ARG:HD3	1:A:447:TYR:HE2	1.71	0.55
1:A:680:TYR:O	1:A:684:VAL:HG23	2.06	0.55
1:A:527:ARG:NH2	1:A:530:ILE:HG12	2.22	0.54
1:A:827:ARG:HH12	1:A:866:ASP:CG	2.16	0.54
1:A:770:GLU:O	1:A:771:LEU:HD23	2.07	0.54
1:A:814:ILE:HG22	1:A:816:LEU:HG	1.89	0.54
1:A:829:TRP:CB	1:A:883:GLY:HA3	2.38	0.54
1:A:740:ILE:HD12	1:A:802:PHE:CE2	2.38	0.54
1:A:483:PHE:O	1:A:492:LEU:HD22	2.07	0.54
1:A:543:ALA:C	1:A:545:LEU:N	2.66	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:625:THR:O	1:A:629:LYS:HB2	2.08	0.54
1:A:822:VAL:HB	1:A:852:VAL:HG22	1.89	0.54
1:A:847:THR:O	1:A:848:ARG:HG3	2.07	0.54
1:A:615:SER:OG	1:A:618:GLU:HB2	2.07	0.54
1:A:688:PHE:CE2	1:A:899:ILE:HD11	2.43	0.54
1:A:860:THR:HG22	1:A:861:LEU:HD23	1.89	0.54
1:A:531:LYS:HG2	1:A:534:ASP:CB	2.34	0.54
1:A:578:VAL:O	1:A:580:GLU:N	2.38	0.53
1:A:791:ILE:HD13	1:A:815:ASN:H	1.73	0.53
1:A:827:ARG:NH2	1:A:866:ASP:HB2	2.23	0.53
1:A:883:GLY:HA2	1:A:886:TRP:NE1	2.18	0.53
1:A:587:ILE:HG22	1:A:588:ALA:N	2.23	0.53
1:A:499:CYS:HB2	1:A:503:VAL:HG23	1.90	0.53
1:A:687:LEU:CD1	1:A:702:MET:HG2	2.39	0.53
1:A:821:ARG:NE	1:A:853:HIS:NE2	2.54	0.53
1:A:440:ILE:CD1	1:A:481:ALA:HA	2.39	0.52
1:A:487:LYS:C	1:A:489:GLU:H	2.16	0.52
1:A:496:LEU:HD12	1:A:497:VAL:H	1.74	0.52
1:A:551:LEU:HG	1:A:556:TRP:HZ2	1.74	0.52
1:A:730:SER:HB3	1:A:733:MET:HB2	1.91	0.52
1:A:527:ARG:O	1:A:528:SER:HB2	2.09	0.52
1:A:506:ASN:O	1:A:510:GLU:HG3	2.10	0.52
1:A:541:THR:HG22	1:A:543:ALA:H	1.75	0.52
1:A:565:GLN:CD	1:A:565:GLN:H	2.18	0.52
1:A:543:ALA:O	1:A:545:LEU:N	2.43	0.51
1:A:514:PHE:C	1:A:516:PRO:HD3	2.35	0.51
1:A:890:LEU:HD12	1:A:895:LEU:HD12	1.93	0.51
1:A:895:LEU:O	1:A:899:ILE:HG12	2.10	0.51
1:A:579:LYS:O	1:A:579:LYS:HG2	2.11	0.51
1:A:628:LYS:C	1:A:630:GLY:H	2.18	0.51
1:A:598:VAL:HG12	1:A:599:ASP:N	2.24	0.51
1:A:680:TYR:CD1	1:A:901:LEU:HB2	2.46	0.51
1:A:540:THR:CG2	1:A:541:THR:N	2.74	0.50
1:A:602:TRP:HB2	1:A:619:PHE:CE2	2.46	0.50
1:A:623:PHE:CD2	1:A:638:LEU:HB2	2.46	0.50
1:A:631:ASP:C	1:A:633:MET:H	2.18	0.50
1:A:441:LYS:O	1:A:442:ALA:HB2	2.11	0.50
1:A:523:PHE:O	1:A:544:VAL:HG11	2.11	0.50
1:A:578:VAL:O	1:A:581:LEU:HG	2.11	0.50
1:A:648:ARG:HH21	1:A:648:ARG:HG3	1.76	0.50
1:A:717:PRO:C	1:A:719:LEU:H	2.19	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:541:THR:CG2	1:A:542:TYR:N	2.75	0.49
1:A:620:LYS:HA	1:A:624:ALA:CB	2.41	0.49
1:A:650:THR:O	1:A:656:ILE:HG21	2.12	0.49
1:A:666:THR:OG1	1:A:853:HIS:ND1	2.45	0.49
1:A:791:ILE:HD13	1:A:814:ILE:HA	1.94	0.49
1:A:468:ALA:HB3	1:A:648:ARG:HB2	1.94	0.49
1:A:527:ARG:NE	1:A:530:ILE:HG12	2.28	0.49
1:A:597:LYS:O	1:A:600:ASP:N	2.45	0.49
1:A:696:GLY:C	1:A:698:LYS:H	2.21	0.49
1:A:740:ILE:HG23	1:A:802:PHE:CE2	2.48	0.49
1:A:459:ASN:HB3	1:A:585:TYR:CD2	2.46	0.49
1:A:571:GLN:HA	1:A:575:PHE:CD1	2.47	0.49
1:A:668:VAL:HG11	1:A:739:ILE:CD1	2.40	0.49
1:A:795:GLN:HE21	1:A:815:ASN:HD21	1.61	0.49
1:A:859:GLY:HA2	1:A:864:LYS:HZ2	1.78	0.49
1:A:435:LEU:HB2	1:A:452:PHE:CE2	2.48	0.48
1:A:505:LYS:HE2	1:A:863:GLU:OE1	2.12	0.48
1:A:776:PRO:HB2	1:A:790:ILE:HG23	1.95	0.48
1:A:474:GLY:O	1:A:478:GLN:HG3	2.13	0.48
1:A:566:ASN:O	1:A:567:ILE:C	2.56	0.48
1:A:438:TYR:HB3	1:A:484:SER:OG	2.13	0.48
1:A:450:LYS:O	1:A:451:GLY:C	2.54	0.48
1:A:688:PHE:C	1:A:690:ASN:N	2.72	0.48
1:A:532:LEU:HD22	1:A:538:ILE:HD13	1.95	0.48
1:A:611:GLY:O	1:A:612:LEU:C	2.57	0.48
1:A:673:THR:H	1:A:676:GLN:HE21	1.62	0.48
1:A:580:GLU:O	1:A:580:GLU:CG	2.62	0.48
1:A:592:THR:C	1:A:594:ILE:H	2.19	0.48
1:A:605:MET:SD	1:A:642:ILE:HG21	2.53	0.48
1:A:623:PHE:C	1:A:626:PRO:HD2	2.39	0.48
1:A:883:GLY:CA	1:A:886:TRP:HE1	2.21	0.48
1:A:777:PHE:O	1:A:790:ILE:HD13	2.14	0.47
1:A:559:ILE:HD12	1:A:583:SER:HB3	1.95	0.47
1:A:597:LYS:NZ	1:A:597:LYS:CB	2.77	0.47
1:A:638:LEU:CG	1:A:642:ILE:HD11	2.40	0.47
1:A:814:ILE:HB	1:A:842:TYR:CD2	2.49	0.47
1:A:706:THR:O	1:A:707:LEU:C	2.57	0.47
1:A:542:TYR:O	1:A:545:LEU:HB3	2.14	0.47
1:A:523:PHE:HB3	1:A:540:THR:HG1	1.77	0.47
1:A:551:LEU:HG	1:A:556:TRP:CZ2	2.49	0.47
1:A:563:GLU:C	1:A:564:ALA:O	2.58	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:605:MET:HA	1:A:605:MET:CE	2.44	0.47
1:A:436:GLU:O	1:A:437:PRO:C	2.57	0.47
1:A:680:TYR:CE2	1:A:710:LEU:HG	2.50	0.47
1:A:433:GLN:O	1:A:435:LEU:N	2.48	0.47
1:A:562:ASP:OD2	3:A:958:PO4:O4	2.32	0.47
1:A:578:VAL:C	1:A:580:GLU:N	2.73	0.47
1:A:698:LYS:O	1:A:698:LYS:HG3	2.14	0.47
1:A:797:ASN:ND2	1:A:798:PRO:HD2	2.25	0.47
1:A:882:SER:O	1:A:886:TRP:CD1	2.67	0.47
1:A:689:ASN:N	1:A:689:ASN:ND2	2.62	0.47
1:A:697:ILE:HG13	1:A:697:ILE:O	2.15	0.47
1:A:864:LYS:HZ2	1:A:902:SER:HB2	1.80	0.47
1:A:653:ASP:HB2	1:A:656:ILE:HG22	1.96	0.46
1:A:785:LYS:O	1:A:789:ASP:CG	2.58	0.46
1:A:500:PRO:O	1:A:501:LEU:C	2.58	0.46
1:A:491:GLU:O	1:A:557:LYS:HD3	2.14	0.46
1:A:584:LYS:HG2	1:A:585:TYR:CE1	2.50	0.46
1:A:827:ARG:NH2	1:A:862:GLU:O	2.49	0.46
1:A:888:THR:OG1	1:A:889:GLU:OE2	2.32	0.46
1:A:594:ILE:HG22	1:A:596:ASN:OD1	2.16	0.46
1:A:593:PRO:O	1:A:594:ILE:C	2.58	0.46
1:A:755:GLN:N	1:A:826:ASP:OD1	2.46	0.46
1:A:759:MET:O	1:A:763:ILE:HG13	2.16	0.46
1:A:624:ALA:O	1:A:625:THR:C	2.58	0.46
1:A:795:GLN:NE2	1:A:815:ASN:ND2	2.64	0.46
1:A:496:LEU:HD12	1:A:497:VAL:N	2.31	0.45
1:A:544:VAL:HG12	1:A:544:VAL:O	2.14	0.45
1:A:840:ARG:NH2	1:A:840:ARG:HG3	2.31	0.45
1:A:771:LEU:O	1:A:772:ASN:O	2.34	0.45
1:A:791:ILE:HD11	1:A:814:ILE:CG2	2.33	0.45
1:A:864:LYS:NZ	1:A:902:SER:HB2	2.32	0.45
1:A:597:LYS:O	1:A:598:VAL:C	2.58	0.45
1:A:728:ARG:CD	1:A:728:ARG:N	2.78	0.45
1:A:620:LYS:HG2	1:A:625:THR:HG23	1.97	0.45
1:A:431:SER:O	1:A:432:PHE:HB2	2.16	0.45
1:A:696:GLY:C	1:A:698:LYS:N	2.75	0.45
1:A:504:LEU:O	1:A:504:LEU:HD23	2.16	0.45
1:A:886:TRP:C	1:A:888:THR:N	2.75	0.45
1:A:891:SER:O	1:A:892:THR:C	2.60	0.45
1:A:516:PRO:C	1:A:518:LEU:N	2.75	0.44
1:A:688:PHE:CE1	1:A:896:ARG:HB2	2.45	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:493:THR:HA	1:A:494:PRO:HA	1.77	0.44
1:A:469:ASP:O	1:A:470:ASP:C	2.58	0.44
1:A:779:TYR:HA	1:A:806:SER:O	2.18	0.44
1:A:814:ILE:O	1:A:842:TYR:HA	2.17	0.44
1:A:894:GLU:O	1:A:897:LYS:HB2	2.17	0.44
1:A:662:ASP:OD1	1:A:845:GLY:O	2.35	0.44
1:A:540:THR:HG23	1:A:541:THR:N	2.33	0.44
1:A:632:ASN:HD22	1:A:635:LYS:HD3	1.83	0.44
1:A:744:LEU:HD11	1:A:802:PHE:CZ	2.53	0.44
1:A:761:LYS:HG3	1:A:777:PHE:HE2	1.82	0.44
1:A:889:GLU:CD	1:A:889:GLU:N	2.72	0.44
1:A:596:ASN:C	1:A:598:VAL:H	2.25	0.44
1:A:648:ARG:O	1:A:648:ARG:HG2	2.17	0.44
1:A:616:TYR:O	1:A:617:SER:C	2.60	0.44
1:A:465:ILE:CA	1:A:608:LEU:HD13	2.48	0.43
1:A:530:ILE:HG23	1:A:532:LEU:HG	1.98	0.43
1:A:586:ARG:HG2	1:A:607:PHE:CZ	2.53	0.43
1:A:900:GLU:HG2	1:A:901:LEU:O	2.18	0.43
1:A:488:LYS:O	1:A:488:LYS:HG3	2.17	0.43
1:A:504:LEU:CD2	1:A:508:GLU:OE1	2.63	0.43
1:A:594:ILE:HG13	1:A:594:ILE:H	1.64	0.43
1:A:450:LYS:O	1:A:453:SER:N	2.51	0.43
1:A:778:LEU:HD23	1:A:805:LEU:HD23	2.00	0.43
1:A:884:ASP:HA	1:A:887:ILE:HD12	2.01	0.43
1:A:811:GLY:O	1:A:812:PHE:C	2.60	0.43
1:A:814:ILE:HB	1:A:842:TYR:HD2	1.83	0.43
1:A:716:HIS:O	1:A:719:LEU:HB2	2.19	0.43
1:A:827:ARG:HD2	1:A:827:ARG:HA	1.70	0.43
1:A:829:TRP:HB2	1:A:883:GLY:HA3	1.99	0.43
1:A:594:ILE:C	1:A:596:ASN:N	2.77	0.43
1:A:698:LYS:O	1:A:698:LYS:CG	2.66	0.43
1:A:527:ARG:NH2	1:A:550:ARG:CB	2.81	0.42
1:A:548:ASP:CG	1:A:549:THR:N	2.74	0.42
1:A:448:GLN:O	1:A:452:PHE:HB3	2.20	0.42
1:A:691:ILE:HG23	1:A:692:ASP:N	2.33	0.42
1:A:564:ALA:C	1:A:566:ASN:N	2.73	0.42
1:A:631:ASP:C	1:A:633:MET:N	2.76	0.42
1:A:860:THR:HG22	1:A:861:LEU:N	2.34	0.42
1:A:525:GLU:HB2	1:A:526:ASP:H	1.60	0.42
1:A:637:GLU:OE1	1:A:637:GLU:HA	2.18	0.42
1:A:441:LYS:N	1:A:441:LYS:CD	2.82	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:487:LYS:C	1:A:489:GLU:N	2.78	0.42
1:A:872:LYS:HG2	1:A:874:SER:OG	2.20	0.42
1:A:7:HIS:O	1:A:8:MET:C	2.63	0.42
1:A:551:LEU:O	1:A:552:LYS:C	2.62	0.42
1:A:605:MET:SD	1:A:642:ILE:CG2	3.08	0.42
1:A:740:ILE:HA	1:A:740:ILE:HD13	1.82	0.42
1:A:499:CYS:HA	1:A:500:PRO:HD3	1.93	0.41
1:A:493:THR:O	1:A:557:LYS:HD2	2.20	0.41
1:A:503:VAL:HA	1:A:506:ASN:OD1	2.20	0.41
1:A:596:ASN:C	1:A:598:VAL:N	2.76	0.41
1:A:840:ARG:HG3	1:A:840:ARG:HH21	1.84	0.41
1:A:869:LEU:HD21	1:A:886:TRP:CZ2	2.54	0.41
1:A:572:THR:C	1:A:574:ILE:N	2.78	0.41
1:A:602:TRP:CD1	1:A:619:PHE:HB2	2.55	0.41
1:A:656:ILE:O	1:A:657:ILE:C	2.63	0.41
1:A:774:GLU:HG2	1:A:776:PRO:HD3	2.02	0.41
1:A:873:ARG:HG3	1:A:876:PHE:CD2	2.43	0.41
1:A:505:LYS:HG2	1:A:669:TYR:CE2	2.55	0.41
1:A:3:PRO:O	1:A:648:ARG:NE	2.54	0.41
1:A:625:THR:HG1	1:A:626:PRO:HD3	1.81	0.41
1:A:555:GLU:HG2	1:A:582:LYS:NZ	2.36	0.41
1:A:682:ALA:C	1:A:684:VAL:H	2.29	0.41
1:A:571:GLN:HA	1:A:575:PHE:CE1	2.56	0.41
1:A:750:ILE:N	1:A:750:ILE:HD12	2.36	0.40
1:A:880:ILE:HG21	1:A:886:TRP:CG	2.56	0.40
1:A:8:MET:O	1:A:9:ALA:CB	2.68	0.40
1:A:559:ILE:O	1:A:559:ILE:HG22	2.21	0.40
1:A:605:MET:HE3	1:A:605:MET:CA	2.49	0.40
1:A:465:ILE:C	1:A:608:LEU:HD13	2.47	0.40
1:A:774:GLU:HG2	1:A:775:VAL:N	2.36	0.40
1:A:551:LEU:O	1:A:554:VAL:N	2.53	0.40
1:A:564:ALA:O	1:A:566:ASN:N	2.54	0.40
1:A:869:LEU:O	1:A:873:ARG:HB2	2.22	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:872:LYS:NZ	1:A:872:LYS:NZ[8_665]	2.15	0.05

5.3 Torsion angles

5.3.1 Protein backbone

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
1	A	476/500 (95%)	350 (74%)	82 (17%)	44 (9%)	0 2

All (44) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	3	PRO
1	A	8	MET
1	A	9	ALA
1	A	525	GLU
1	A	526	ASP
1	A	564	ALA
1	A	567	ILE
1	A	572	THR
1	A	879	ILE
1	A	885	SER
1	A	892	THR
1	A	527	ARG
1	A	544	VAL
1	A	546	LEU
1	A	570	PRO
1	A	571	GLN
1	A	579	LYS
1	A	772	ASN
1	A	780	GLY
1	A	880	ILE
1	A	438	TYR
1	A	517	HIS
1	A	565	GLN
1	A	568	LYS
1	A	593	PRO
1	A	612	LEU
1	A	624	ALA
1	A	689	ASN

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Mol	Chain	Res	Type
1	A	770	GLU
1	A	12	LYS
1	A	524	HIS
1	A	549	THR
1	A	683	GLU
1	A	693	SER
1	A	694	VAL
1	A	782	LEU
1	A	811	GLY
1	A	7	HIS
1	A	591	GLY
1	A	598	VAL
1	A	619	PHE
1	A	781	GLU
1	A	594	ILE
1	A	626	PRO

5.3.2 Protein sidechains ⓘ

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	419/446 (94%)	391 (93%)	28 (7%)	15	46

All (28) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	452	PHE
1	A	456	ARG
1	A	461	LEU
1	A	501	LEU
1	A	502	SER
1	A	511	LEU
1	A	524	HIS
1	A	538	ILE
1	A	540	THR
1	A	551	LEU

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Mol	Chain	Res	Type
1	A	597	LYS
1	A	604	ILE
1	A	632	ASN
1	A	642	ILE
1	A	647	LEU
1	A	665	GLU
1	A	670	CYS
1	A	671	ASN
1	A	728	ARG
1	A	733	MET
1	A	735	ARG
1	A	738	GLU
1	A	740	ILE
1	A	752	ILE
1	A	791	ILE
1	A	803	ILE
1	A	828	TRP
1	A	867	GLN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (11) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	478	GLN
1	A	565	GLN
1	A	571	GLN
1	A	632	ASN
1	A	676	GLN
1	A	689	ASN
1	A	712	GLN
1	A	755	GLN
1	A	795	GLN
1	A	797	ASN
1	A	836	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

Of 11 ligands modelled in this entry, 3 are monoatomic - leaving 8 for Mogul analysis.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
3	PO4	A	956	-	4,4,4	1.71	0	6,6,6	0.48	0
3	PO4	A	959	-	4,4,4	1.69	0	6,6,6	0.46	0
3	PO4	A	958	-	4,4,4	1.76	2 (50%)	6,6,6	0.44	0
3	PO4	A	957	-	4,4,4	1.63	1 (25%)	6,6,6	0.50	0
3	PO4	A	960	-	4,4,4	1.73	1 (25%)	6,6,6	0.44	0
3	PO4	A	954	-	4,4,4	1.71	0	6,6,6	0.47	0
3	PO4	A	955	-	4,4,4	1.72	0	6,6,6	0.48	0
3	PO4	A	961	-	4,4,4	1.71	0	6,6,6	0.46	0

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	958	PO4	P-O4	-2.15	1.48	1.54
3	A	957	PO4	P-O2	-2.09	1.48	1.54
3	A	958	PO4	P-O3	-2.01	1.48	1.54
3	A	960	PO4	P-O3	-2.00	1.48	1.54

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	958	PO4	1	0
3	A	961	PO4	1	0

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

EDS was not executed - this section is therefore empty.

6.2 Non-standard residues in protein, DNA, RNA chains

EDS was not executed - this section is therefore empty.

6.3 Carbohydrates

EDS was not executed - this section is therefore empty.

6.4 Ligands

EDS was not executed - this section is therefore empty.

6.5 Other polymers

EDS was not executed - this section is therefore empty.